

# Beamline 17-BM / IMCA-CAT

**Scientific focus:** Pharmaceutical macromolecular crystallography

**Scientific programs:** Structures of protein–ligand complexes, *de novo* protein structures, drug design, protein engineering, and crystallographic methods development

## Optics & Optical Performance

- Daresbury double-crystal constant off-set monochromator
  - 28 m from source
  - 6–20 keV energy range Si(111)
  - $10^{-4}$  energy resolution ( $\Delta E/E$ ) at 10 keV
  - 35 mm offset below orbital plane
  - water cooling
- sagittally bent 2nd monochromator crystal
- Daresbury vertically focusing mirror
- $3 \times 10^{11}$  ph/sec flux on  $200 \mu\text{m} \times 200 \mu\text{m}$  sample at 12.4 keV

## Experiment Stations

### 17-BM-A

- white beam first optics enclosure

### 17-BM-B

- white/monochromatic beam station
- monochromatic macromolecular crystallography

## Detectors

- Mar 165 CCD
- fluorescence detector

## Beamline Controls and Data Acquisition

- controls: Sun and Linux systems running EPICS with VME “MX” software (locally developed), running on UNIX
- data acquisition: proprietary software from Mar

## Beamline Support Equipment/Facilities

- 4° chill room in wet lab
- user-accessible computers for data processing
- Oxford Instruments cryojet for sample cooling

## Bending Magnet Source Characteristics (nominal)

source	APS bending magnet
critical energy	19.51 keV
on-axis peak brilliance at 16.3 keV	$2.9 \times 10^{15}$ ph/sec/mrad <sup>2</sup> /mm <sup>2</sup> /0.1%bw
on-axis peak angular flux at 16.3 keV	$9.6 \times 10^{13}$ ph/sec/mrad <sup>2</sup> /0.1%bw
on-axis peak horizontal angular flux at 5.6 keV	$1.6 \times 10^{13}$ ph/sec/mradh/0.1%bw
source size at critical energy $\sum_x$ $\sum_y$	145 $\mu\text{m}$ 36 $\mu\text{m}$
source divergence at critical energy $\sum_{x'}$ $\sum_{y'}$	6 mrad 47 $\mu\text{rad}$